

chain nodes :

15

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13

chain bonds :

12-15

ring bonds :

1-2 1-6 2-3 3-4 4-5 4-7 5-6 5-10 7-8 7-11 8-9 8-13 9-10 11-12 12-13

exact/norm bonds :

8-13 12-13 12-15

exact bonds :

4-7 5-10 7-8 7-11 8-9 9-10 11-12

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6

isolated ring systems :

containing 1 :

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom  
12:Atom 13:Atom 15:CLASS

Trying 3106016892...Open

```
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NEWS	3	JUN 20 WIPO/PCT Patents Fulltext Database now on STN
NEWS	4	JUN 28 CAS covers Web-distributed preprints
NEWS	5	JUL 7 Patent Full-text Cluster, PNTTEXT, introduced
NEWS	6	JUL 27 EUROPATFULL - loading of backlog data
NEWS	7	JUL 27 MORE FREQUENT UPDATES FOR DERWENT WORLD PATENTS INDEX IN 2000
NEWS	8	JUL 27 Derwent Journal Of Synthetic Methods Reloaded with New Data
NEWS	9	JUL 27 DERWENT WORLD PATENTS INDEX: FAST TRACK RELEASE OF EQUIVALENT PATENTS
NEWS	10	AUG 21 Instant Access to FDA Regulatory Information with DIOGENES
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ENTRY

TOTAL  
SESSION

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Page 1

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0.15

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DICTIONARY FILE UPDATES: 8 SEP 2000 HIGHEST RN 288568-75-2

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=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>  
Uploading c:\stnexp4\queries\09552969.str

L1 STRUCTURE uploaded

=> que L1

L2 QUE L1

=> s 11

SAMPLE SEARCH INITIATED 12:29:24 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 159 TO ITERATE

100.0% PROCESSED 159 ITERATIONS 0 ANSWERS  
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 2424 TO 3936  
PROJECTED ANSWERS: 0 TO 0

L3 0 SEA SSS SAM L1

=> s 11 ful

FULL SEARCH INITIATED 12:29:37 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 3220 TO ITERATE

100.0% PROCESSED 3220 ITERATIONS 12 ANSWERS  
SEARCH TIME: 00.00.01

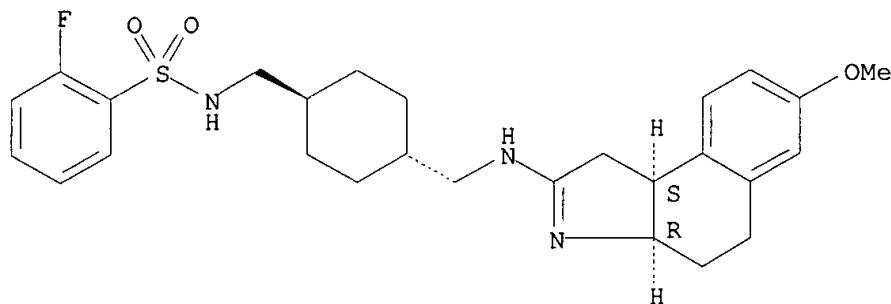
L4 12 SEA SSS FUL L1

=> d scan

L4 12 ANSWERS REGISTRY COPYRIGHT 2000 ACS  
IN Benzenesulfonamide,  
2-fluoro-N-[{trans-4-[[[(3aR,9bS)-3a,4,5,9b-tetrahydro-  
7-methoxy-1H-benz[e]indol-2-yl]amino]methyl]cyclohexyl}methyl]-,

monohydrochloride, rel- (9CI)  
MF C27 H34 F N3 O3 S . Cl H

Relative stereochemistry.



• HC1

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> fil caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	126.60	126.75

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FILE COVERS 1967 - 9 Sep 2000 VOL 133 ISS 12  
FILE LAST UPDATED: 8 Sep 2000 (20000908/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

This file supports REG1stRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

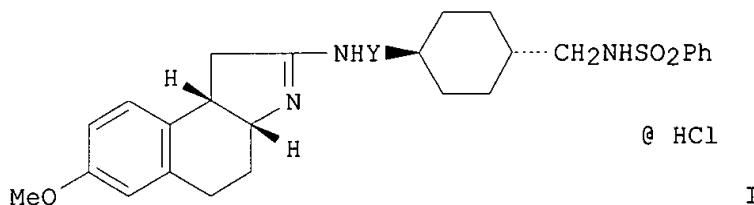
Now you can extend your author, patent assignee, patent information, and title searches back to 1907. The records from 1907-1966 now have this searchable data in CAOLD. You now have electronic access to all of CA: 1907 to 1966 in CAOLD and 1967 to the present in CAPLUS on STN.

=> s 14

L5 1 L4

=> d ibib abs hitstr 1-

L5 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2000 ACS  
 ACCESSION NUMBER: 2000:146875 CAPLUS  
 DOCUMENT NUMBER: 132:279085  
 TITLE: N-(Sulfonamidoalkyl)(tetrahydro-1H-benz[e]indol-2-yl)amines: potent antagonists of human neuropeptide Y Y5 receptor  
 AUTHOR(S): McNally, James J.; Youngman, Mark A.; Lovenberg, Timothy W.; Nepomuceno, Diane H.; Wilson, Sandy J.; Dax, Scott L.  
 CORPORATE SOURCE: Drug Discovery, The R. W. Johnson Pharmaceutical Research Institute, Spring House, PA, 19477, USA  
 SOURCE: Bioorg. Med. Chem. Lett. (2000), 10(3), 213-216  
 PUBLISHER: Elsevier Science Ltd. Feb. 7, 2000  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 GI



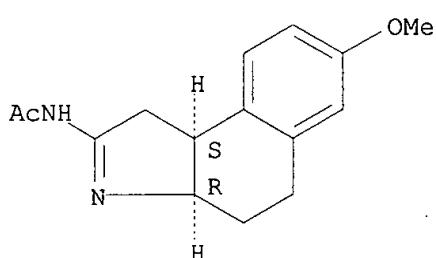
AB (3A,4,5,9b-Tetrahydro-1H-benz[e]indol-2-yl)amines were prep'd. via reductive amination and concomitant cyclization of .alpha.-cyanomethyl-.beta.-aminotetralins. N-acylation with .OMEGA.-sulfonamido-carboxylic acids and subsequent redn. afforded a series of title compds., e.g., I (Y = CO, CH<sub>2</sub>), which bound to the human neuropeptide Y Y5 receptor with nanomolar affinity.

IT 263843-82-9P 263843-83-0P 263843-84-1P  
 263843-85-2P 263843-86-3P 263843-87-4P  
 263843-88-5P 263843-89-6P 263843-90-9P  
 263843-91-0P 263843-92-1P 263843-93-2P  
 RL: BAC (Biological activity or effector, except adverse); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation)  
 (prepn. and neuropeptide Y5 receptor binding affinity of)

RN 263843-82-9 CAPLUS

CN Acetamide, N-[(3aR,9bS)-3a,4,5,9b-tetrahydro-7-methoxy-1H-benz[e]indol-2-yl]-, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

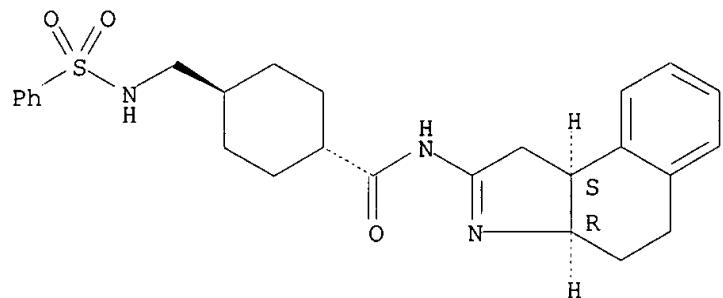


RN 263843-83-0 CAPLUS

CN Cyclohexanecarboxamide, 4-[[ (phenylsulfonyl)amino]methyl]-N-[(3aR,9bS)-

3a,4,5,9b-tetrahydro-1H-benz[e]indol-2-yl]-, monohydrochloride,  
trans-rel-  
(9CI) (CA INDEX NAME)

Relative stereochemistry.

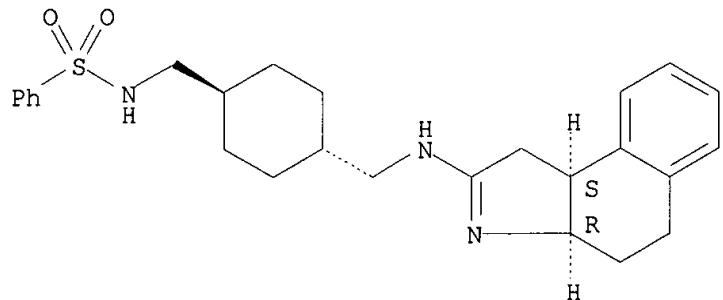


● HCl

RN 263843-84-1 CAPLUS

CN Benzenesulfonamide, N-[{trans-4-[[[(3aR,9bS)-3a,4,5,9b-tetrahydro-1H-benz[e]indol-2-yl]amino]methyl]cyclohexyl}methyl]-, monohydrochloride, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

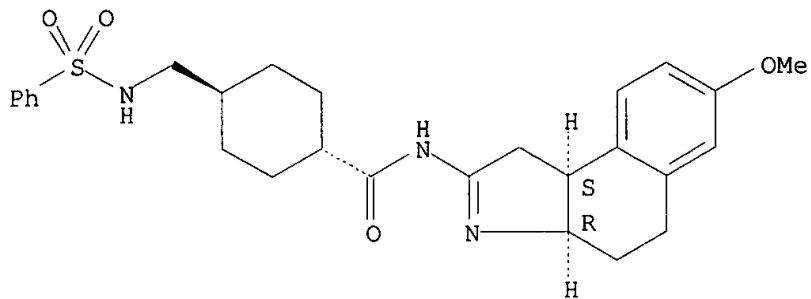


● HCl

RN 263843-85-2 CAPLUS

CN Cyclohexanecarboxamide, 4-[(phenylsulfonyl)amino]methyl]-N-[(3aR,9bS)-3a,4,5,9b-tetrahydro-7-methoxy-1H-benz[e]indol-2-yl]-, monohydrochloride, trans-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

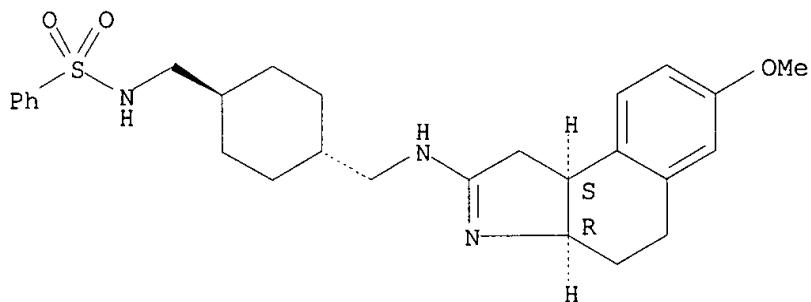


• HCl

RN 263843-86-3 CAPLUS

CN Benzenesulfonamide, N-[{trans-4-[[[(3aR,9bS)-3a,4,5,9b-tetrahydro-7-methoxy-1H-benz[e]indol-2-yl]amino]methyl]cyclohexyl}methyl]-, monohydrochloride, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

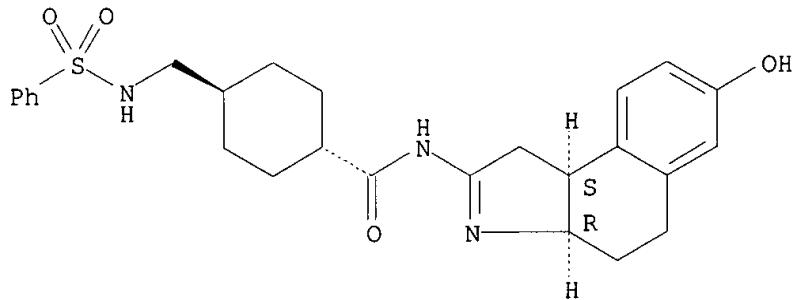


• HCl

RN 263843-87-4 CAPLUS

CN Cyclohexanecarboxamide, 4-[[{(phenylsulfonyl)amino}methyl]-N-[(3aR,9bS)-3a,4,5,9b-tetrahydro-7-hydroxy-1H-benz[e]indol-2-yl]-, monohydrochloride, trans-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

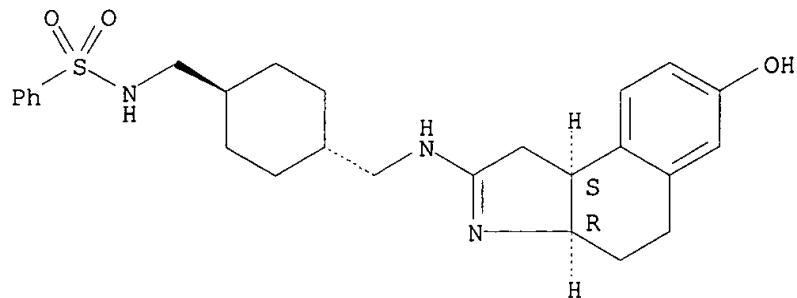


• HCl

RN 263843-88-5 CAPLUS

CN Benzenesulfonamide, N-[[trans-4-[[[(3aR,9bS)-3a,4,5,9b-tetrahydro-7-hydroxy-1H-benz[e]indol-2-yl]amino]methyl]cyclohexyl]methyl]-, monohydrochloride, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

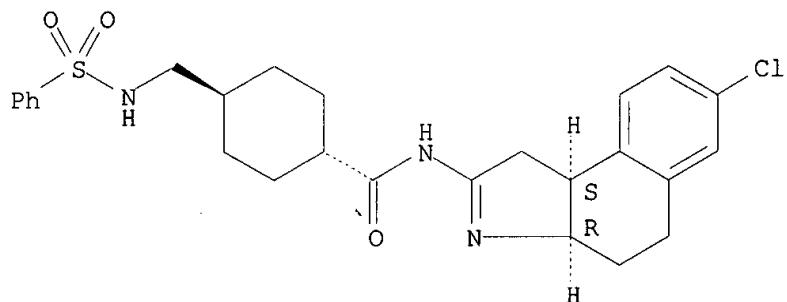


• HCl

RN 263843-89-6 CAPLUS

CN Cyclohexanecarboxamide, N-[(3aR,9bS)-7-chloro-3a,4,5,9b-tetrahydro-1H-benz[e]indol-2-yl]-4-[(phenylsulfonyl)amino]methyl-, monohydrochloride, trans-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

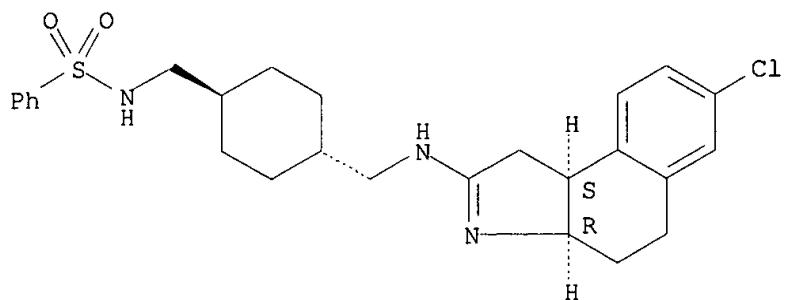


• HCl

RN 263843-90-9 CAPLUS

CN Benzenesulfonamide,  
N-[[trans-4-[[[(3aR,9bS)-7-chloro-3a,4,5,9b-tetrahydro-1H-benz[e]indol-2-yl]amino]methyl]cyclohexyl]methyl]-, monohydrochloride, rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

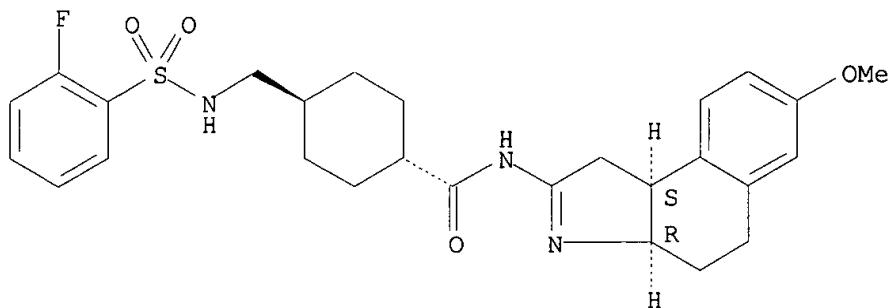


• HCl

RN 263843-91-0 CAPLUS

CN Cyclohexanecarboxamide, 4-[[[(2-fluorophenyl)sulfonyl]amino]methyl]-N-[(3aR,9bS)-3a,4,5,9b-tetrahydro-7-methoxy-1H-benz[e]indol-2-yl]-, monohydrochloride, trans-rel- (9CI) (CA INDEX NAME)

Relative stereochemistry.

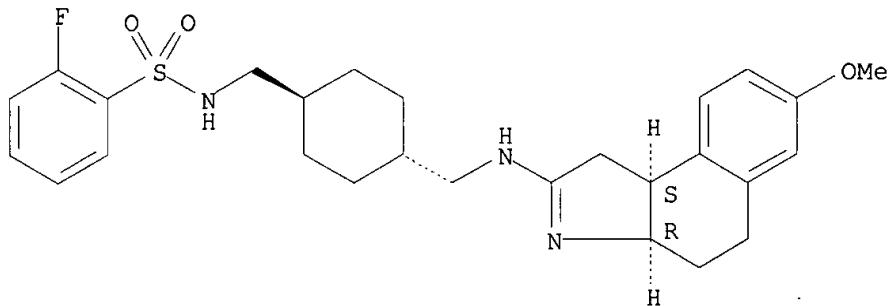


• HCl

RN 263843-92-1 CAPLUS

CN Benzenesulfonamide, 2-fluoro-N-[(trans-4-[[[(3aR,9bS)-3a,4,5,9b-tetrahydro-7-methoxy-1H-benz[e]indol-2-yl]amino]methyl]cyclohexyl)methyl]-, monohydrochloride, rel- (9CI) (CA INDEX NAME)

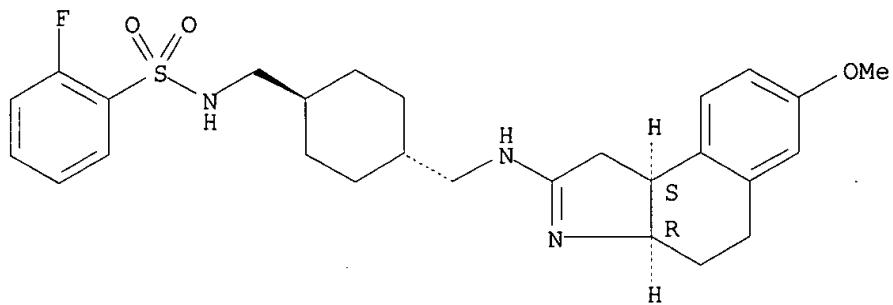
Relative stereochemistry.



• HCl

Hong

Page 8

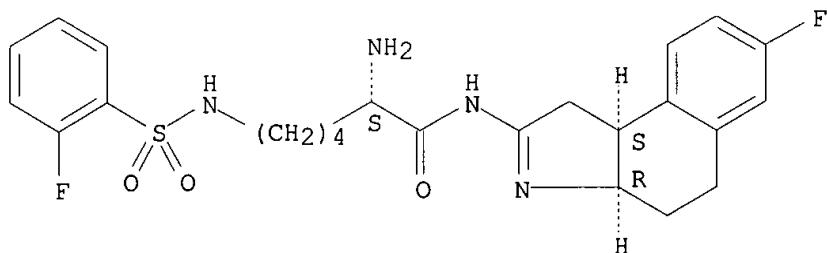


• HCl

RN 263843-93-2 CAPLUS

CN Hexanamide, 2-amino-6-[(2-fluorophenyl)sulfonyl]amino]-N-[ (3aR,9bS)-7-fluoro-3a,4,5,9b-tetrahydro-1H-benz[e]indol-2-yl]-, monohydrochloride, (2S)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.



• HCl

REFERENCE COUNT:

20

REFERENCE(S):

- (1) Castan, I; Endocrinology 1992, V131, P1970 CAPLUS
- (2) Clark, J; Endocrinology 1984, V115, P427 CAPLUS
- (3) Criscione, L; J Clin Invest 1998, V102, P2136 CAPLUS
- (4) Fuhlendorff, J; Proc Natl Acad Sci USA 1990, V87, P182 CAPLUS
- (5) Gerald, C; WO 97/46250 1997 CAPLUS

ALL CITATIONS AVAILABLE IN THE RE FORMAT

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COST IN U.S. DOLLARS

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4.90 131.65

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=>  
Uploading c:\stnexp4\queries\09552969.str

L6 STRUCTURE uploaded

=> que L6

L7 QUE L6

=> s 16

SAMPLE SEARCH INITIATED 12:32:54 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 179 TO ITERATE

100.0% PROCESSED 179 ITERATIONS 0 ANSWERS  
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 2778 TO 4382  
PROJECTED ANSWERS: 0 TO 0

L8 0 SEA SSS SAM L6

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>  
Uploading c:\stnexp4\queries\09552969.str

L9 STRUCTURE uploaded

=> que L9

L10 QUE L9

=> s 19

SAMPLE SEARCH INITIATED 12:33:43 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 19192 TO ITERATE

5.2% PROCESSED 1000 ITERATIONS 10 ANSWERS  
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)  
SEARCH TIME: 00.00.01

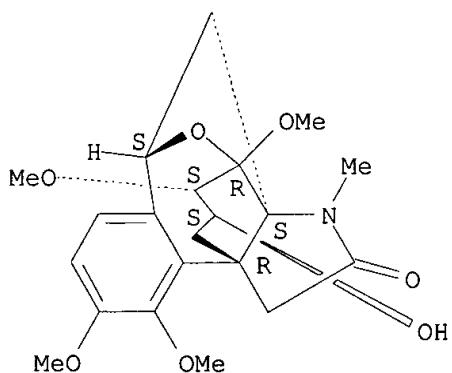
FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 375589 TO 392091  
PROJECTED ANSWERS: 3007 TO 4669

L11 10 SEA SSS SAM L9

=> d scan

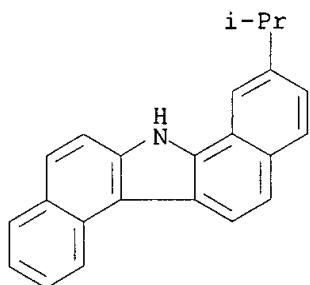
L11 10 ANSWERS REGISTRY COPYRIGHT 2000 ACS  
IN Hasubanan-16-one, 8,10-epoxy-6-hydroxy-3,4,7,8-tetramethoxy-17-methyl-,  
(6. $\beta$ .,7. $\beta$ .,8. $\beta$ .,10. $\beta$ .)- (9CI)  
MF C21 H27 N O7

Absolute stereochemistry.



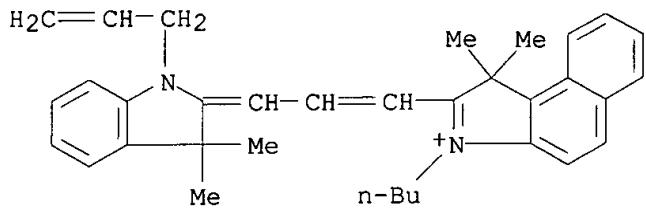
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):2

L11 10 ANSWERS REGISTRY COPYRIGHT 2000 ACS  
IN 7H-Dibenzo[a,g]carbazole, 9-isopropyl- (8CI)  
MF C23 H19 N

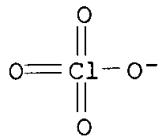


L11 10 ANSWERS REGISTRY COPYRIGHT 2000 ACS  
IN 1H-Benz[e]indolium,  
3-butyl-2-[3-[1,3-dihydro-3,3-dimethyl-1-(2-propenyl)-  
2H-indol-2-ylidene]-1-propenyl]-1,1-dimethyl-, perchlorate (9CI)  
MF C34 H39 N2 . Cl O4

CM 1



CM 2



HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>  
Uploading c:\stnexp4\queries\09552969.str

L12 STRUCTURE UPLOADED

=> que L12

L13 QUE L12

=> s l12

SAMPLE SEARCH INITIATED 12:34:59 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 19192 TO ITERATE

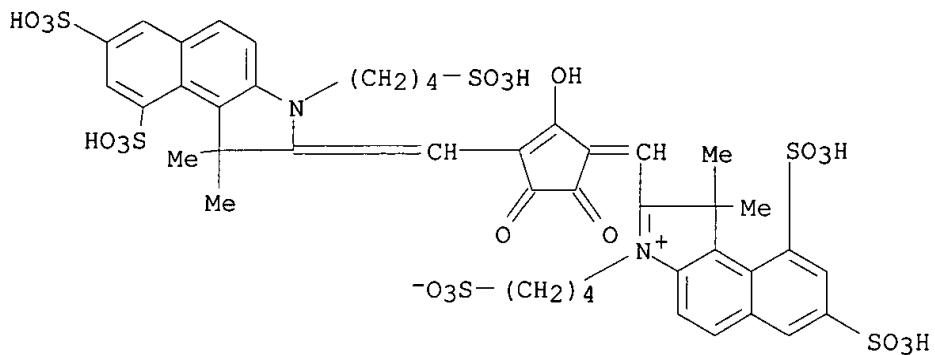
5.2% PROCESSED 1000 ITERATIONS 10 ANSWERS  
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)  
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 375589 TO 392091  
PROJECTED ANSWERS: 3007 TO 4669

L14 10 SEA SSS SAM L12

=> d scan

L14 10 ANSWERS REGISTRY COPYRIGHT 2000 ACS  
IN 1H-Benz[e]indolium, 2-[3-[(1,3-dihydro-1,1-dimethyl-7,9-disulfo-3-(4-sulfobutyl)-2H-benz[e]indol-2-ylidene)methyl]-2-hydroxy-4,5-dioxo-2-cyclopenten-1-ylidene)methyl]-1,1-dimethyl-7,9-disulfo-3-(4-sulfobutyl)-, inner salt, hexapotassium salt (9CI)  
MF C43 H44 N2 O21 S6 . 6 K

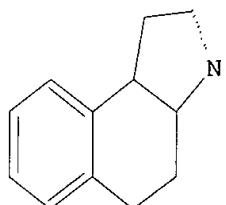


• 6 K

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> d 112

L12 HAS NO ANSWERS  
L12 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 112 ful

FULL SEARCH INITIATED 12:35:42 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 387071 TO ITERATE

100.0% PROCESSED 387071 ITERATIONS  
SEARCH TIME: 00.00.07

4659 ANSWERS

L15 4659 SEA SSS FUL L12

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>  
Uploading c:\stnexp4\queries\09552969.str

L16 STRUCTURE UPLOADED

=> que L16

L17 QUE L16

=> s 116 ful sub=115

FULL SUBSET SEARCH INITIATED 12:36:58 FILE 'REGISTRY'  
FULL SUBSET SCREEN SEARCH COMPLETED - 14 TO ITERATE

100.0% PROCESSED 14 ITERATIONS 12 ANSWERS  
SEARCH TIME: 00.00.01

L18 12 SEA SUB=L15 SSS FUL L16

=> fil caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	158.40	290.05
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.56

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FILE LAST UPDATED: 8 Sep 2000 (20000908/ED)

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=> s l18

L19 1 L18

=> d ibib

L19 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2000 ACS  
ACCESSION NUMBER: 2000:146875 CAPLUS  
DOCUMENT NUMBER: 132:279085  
TITLE: N-(Sulfonamidoalkyl)(tetrahydro-1H-benz[e]indol-2-yl)amines: potent antagonists of human neuropeptide Y Y5 receptor

AUTHOR(S): McNally, James J.; Youngman, Mark A.; Lovenberg,  
 Timothy W.; Nepomuceno, Diane H.; Wilson, Sandy J.;  
 Dax, Scott L.  
 CORPORATE SOURCE: Drug Discovery, The R. W. Johnson Pharmaceutical  
 Research Institute, Spring House, PA, 19477, USA  
 SOURCE: Bioorg. Med. Chem. Lett. (2000), 10(3), 213-216  
 CODEN: BMCLE8; ISSN: 0960-894X  
 PUBLISHER: Elsevier Science Ltd.  
 DOCUMENT TYPE: Journal  
 LANGUAGE: English  
 REFERENCE COUNT: 20  
 REFERENCE(S):  
     (1) Castan, I; Endocrinology 1992, V131, P1970 CAPLUS  
     (2) Clark, J; Endocrinology 1984, V115, P427 CAPLUS  
     (3) Criscione, L; J Clin Invest 1998, V102, P2136  
         CAPLUS  
     (4) Fuhlendorff, J; Proc Natl Acad Sci USA 1990, V87,  
         P182 CAPLUS  
     (5) Gerald, C; WO 97/46250 1997 CAPLUS  
 ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> fil reg

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	1.49	291.54
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
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FILE 'REGISTRY' ENTERED AT 12:38:24 ON 09 SEP 2000  
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STRUCTURE FILE UPDATES: 8 SEP 2000 HIGHEST RN 288568-75-2  
 DICTIONARY FILE UPDATES: 8 SEP 2000 HIGHEST RN 288568-75-2

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 11, 2000

Please note that search-term pricing does apply when  
 conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT  
 for details.

=> d his

(FILE 'HOME' ENTERED AT 12:29:03 ON 09 SEP 2000)

FILE 'REGISTRY' ENTERED AT 12:29:06 ON 09 SEP 2000

L1                   STRUCTURE UPLOADED  
 L2                   QUE L1  
 L3                   0 S L1  
 L4                   12 S L1 FUL

FILE 'CAPLUS' ENTERED AT 12:30:35 ON 09 SEP 2000

L5                   1 S L4

FILE 'REGISTRY' ENTERED AT 12:32:36 ON 09 SEP 2000

L6                   STRUCTURE UPLOADED  
 L7                   QUE L6

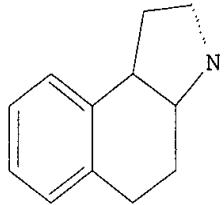
```
L8      0 S L6
L9      STRUCTURE uploaded
L10     QUE L9
L11     10 S L9
L12      STRUCTURE uploaded
L13     QUE L12
L14     10 S L12
L15     4659 S L12 FUL
L16      STRUCTURE uploaded
L17     QUE L16
L18     12 S L16 FUL SUB=L15
```

```
FILE 'CAPLUS' ENTERED AT 12:37:11 ON 09 SEP 2000
L19      1 S L18
```

```
FILE 'REGISTRY' ENTERED AT 12:38:24 ON 09 SEP 2000
```

```
=> d 112
```

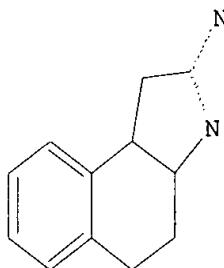
```
L12 HAS NO ANSWERS
L12      STR
```



Structure attributes must be viewed using STN Express query preparation.

```
=> d 116
```

```
L16 HAS NO ANSWERS
L16      STR
```



Structure attributes must be viewed using STN Express query preparation.

```
=> save 112
```

```
ENTER NAME OR (END):f9552969/a
```

```
L12 IS NOT AN ANSWER SET
```

```
A name ending in /A can only be used to save the L-number answer
set resulting from a search.
```

=> save l15

ENTER NAME OR (END) :f9552969/a

ANSWER SET L15 HAS BEEN SAVED AS 'F9552969/A'

=> logoff

ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF  
LOGOFF? (Y)/N/HOLD:y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.90	292.44
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.56

STN INTERNATIONAL LOGOFF AT 12:40:20 ON 09 SEP 2000

Trying 3106016892...Open

```
Welcome to STN International! Enter x:x  
LOGINID:ssspta1611hx1  
PASSWORD:  
TERMINAL (ENTER 1, 2, 3, OR ?):2
```

\* \* \* \* \* \* \* \* \* \* \* \* \* \* \* Welcome to STN International \* \* \* \* \* \* \* \* \* \* \*

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NEWS 3 Jun 20 WIPO/PCT Patents Fulltext Database now on STN

NEWS 4 Jun 28 CAS covers Web-distributed preprints

NEWS 5 Jul 7 Patent Full-text Cluster, PNTTEXT, introduced

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FILE 'HOME' ENTERED AT 10:35:25 ON 18 SEP 2000

=> activate

THIS COMMAND NOT AVAILABLE IN THE CURRENT FILE  
Some commands only work in certain files. For example, the EXPAND command can only be used to look at the index in a file which has an index. Enter "HELP COMMANDS" at an arrow prompt (=>) for a list of commands which can be used in this file.

=> fil reg

COST IN U.S. DOLLARS	SINCE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.15	0.15

FILE 'REGISTRY' ENTERED AT 10:35:42 ON 18 SEP 2000  
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STRUCTURE FILE UPDATES: 17 SEP 2000 HIGHEST RN 289617-02-3  
DICTIONARY FILE UPDATES: 17 SEP 2000 HIGHEST RN 289617-02-3

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 11, 2000

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Structure search limits have been increased. See HELP SLIMIT for details.

=> activate

ENTER NAME OF SAVED ITEM TO ACTIVATE OR (END):f9552969

'F9552969' MUST END IN '/Q', '/A', '/L', '/S', OR '/B'  
The name of a saved query (or structure or screen set) ends in '/Q'.  
The name of a saved answer set ends in '/A'. The name of a saved L# list ends in '/L'. The name of an SDI request ends in '/S'. The name of a BATCH search request ends in '/B'. You must enter the /Q, /A, /L, /S, or /B at the end of the name.  
ENTER NAME OF SAVED ITEM TO ACTIVATE OR (END):f9552969/a

L1 STR  
L2 4659 SEA FILE=REGISTRY SSS FUL L1

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>  
Uploading c:\stnexp4\queries\09552969.str

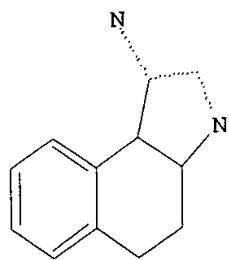
L3 STRUCTURE uploaded

=> que L3

L4 QUE L3

=> d 13

L3 HAS NO ANSWERS  
L3 STR



Structure attributes must be viewed using STN Express query preparation.

=> s 13

SAMPLE SEARCH INITIATED 10:40:05 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 6 TO ITERATE

100.0% PROCESSED 6 ITERATIONS 0 ANSWERS  
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 6 TO 266  
PROJECTED ANSWERS: 0 TO 0

L5 0 SEA SSS SAM L3

=> s 13 ful

FULL SEARCH INITIATED 10:40:18 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 121 TO ITERATE

100.0% PROCESSED 121 ITERATIONS 0 ANSWERS  
SEARCH TIME: 00.00.01

L6 0 SEA SSS FUL L3

=> logoff

ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF  
LOGOFF? (Y)/N/HOLD:y

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	128.40	128.55

STN INTERNATIONAL LOGOFF AT 10:40:43 ON 18 SEP 2000